

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
 Data File : VN070082.D
 Acq On : 14 Dec 2021 20:06
 Operator : JC/MD
 Sample : M5017-49ME
 Misc : 5.08g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP1-BME

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/15/2021
 Supervised By : Mahesh Dadoda 12/21/2021

Quant Time: Dec 15 05:22:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.080	168	1164385	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	2251890	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2589890	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1191929	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.432	65	951268	52.913	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 105.820%			
35) Dibromofluoromethane	8.019	113	627518	44.488	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 88.980%			
50) Toluene-d8	10.440	98	3025738	53.423	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.840%			
62) 4-Bromofluorobenzene	12.731	95	1294666	63.097	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 126.200%			
Target Compounds						
11) Tert butyl alcohol	5.294	59	34032	10.014	ug/l #	74
16) Acetone	4.304	43	15995m	1.479	ug/l	
18) Methyl Acetate	4.867	43	100373	3.075	ug/l #	81
25) 2-Butanone	7.345	43	110453	7.517	ug/l	92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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